

Light Glueball masses using the Multilevel Algorithm.

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Following the multilevel scheme we present an error reduction algorithm for extracting glueball masses from monte-carlo simulations of pure SU(3) lattice gauge theory. We look at the two lightest states viz. the 0^{++} and 2^{++} . Our method involves looking at correlations between large wilson loops and does not require any smearing of links. The error bars we obtain are at the moment comparable to those obtained using smeared operators. We also present a comparison of our method with the naive method.

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1. Introduction

Low lying spectrum of pure Yang-Mills theory consists of glueballs. Quarks allow glueballs to decay but their signature is expected to remain in the QCD spectrum. While glueballs have not yet been discovered experimentally, there are candidates [1]. For a recent review on the status of glueballs see reference [2]. Glueball correlators can be computed in lattice gauge theory simulations but extraction of glueball masses from correlation functions are extremely difficult because the correlation functions are dominated by statistical noise. Recent attempts to compute glueball masses in both pure Yang-Mills theory and lattice QCD using different strategies to combat the statistical noise have been reported in references [3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15, 16, 17]. In this article we report on our attempts to reduce the statistical noise on glueball correlators obtained from simulations of pure SU(3) Yang-Mills theory in 4 Euclidean dimensions.

We explored the scalar (0^{++}) and the tensor (2^{++}) channels and to reduce the noise on the correlators in these channels, we tried the following two strategies (i) construct glueball operators from large wilson loops with extents of about half a fermi in each direction and (ii) extract masses from the correlators with fit range as large as possible (0.5 – 1.0 fermi) to reduce contamination from excited states. The idea to use large Wilson loops to construct glueball operators was proposed in [18] but it was only recently that it was coupled to error reduction techniques to estimate the expectation values of the Wilson loops accurately [19].

2. Algorithm

For updating the pure Yang-Mills fields we used the Cabibo-Marinari heatbath for SU(3) with 3 over-relaxation steps for every heatbath step. Between each measurement we did 10 full sweeps of the lattice to ensure that successive measurements could be treated as independent.

The glueball operators were constructed using Wilson loops. For the scalar channel we constructed the temporal correlator between the operators ($\mathcal{A} - \langle \mathcal{A} \rangle$) at different time slices with $\mathcal{A} = \text{Re}(P_{xy} + P_{xz} + P_{yz})$ and for the tensor channel we took the two operators $\mathcal{E}_1 = \text{Re}(P_{xz} - P_{yz})$ and $\mathcal{E}_2 = \text{Re}(P_{xz} + P_{yz} - 2P_{xy})$. Here P_{ij} denotes a Wilson loop in the ij plane with ij going over the spatial directions.

For the noise reduction scheme we used the philosophy of multilevel algorithm [20]. This method is particularly useful in theories with a mass gap, where the distant regions of the theory are uncorrelated as the correlation length is finite.

The principle of multilevel algorithm is to compute the expectation values in a nested manner. Intermediate values are first constructed by averaging over sub-lattices with boundaries and then the full expectation values are obtained by averging over the intermediate values with different boundaries obtained by updating the full lattice. The intermediate averages can be computed in a nested manner and our innermost noise reduction step was to use a semi-analytic multihit on the SU(3) links [21] with which the Wilson loops were constructed. This reduced the fluctuations in the expectation values of the glueball operators. The multilevel, on top of the multihit, was used to reduce the fluctuation in the correlators.

In figure 1 we illustrate our slicing of the lattice and the computation of the intermediate expectation values of glueball operators (Wilson loops) by performing several sub-lattice updates.

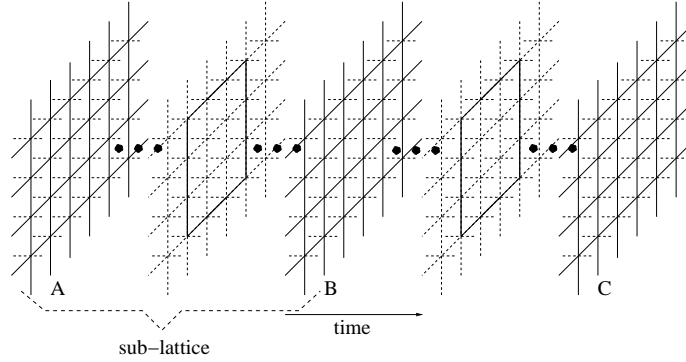


Figure 1: Slicing the lattice for the multilevel update. Solid links on slices A,B and C are frozen during the sub-lattice updates. Thick lines in the Wilson loops indicate that the corresponding links have been replaced by their multihit averages.

Lattice Size	β	(r_0/a_t)	sub-lattice thickness	iupd	loop size
$10^3 \times 18$	5.7	2.922(9)	3	30	2×2
$12^3 \times 18$	5.8	3.673(5)	3	25	3×3
$16^3 \times 24$	5.95	4.898(12)	4	50	5×5

Table 1: Simulation parameters for the scalar channel

Lattice Size	β	(r_0/a_t)	sub-lattice thickness	iupd	loop size
$12^3 \times 18$	5.8	3.673(5)	3	70	3×3
$12^3 \times 20$	5.95	4.898(12)	5	100	5×5
$12^3 \times 20$	6.07	6.033(17)	5	100	5×5

Table 2: Simulation parameters for the tensor channel

The multilevel algorithm has a number of extra parameters such as the thickness of the sub-lattice and the number of sublattice updates used in the intermediate averages. Optimal values of these parameters depend on the observable one is trying to estimate and some tuning of these parameters are essential for efficient error reduction. In tables 1 and 2 we record these along with other simulation parameters for our runs. The scale is set through the Sommer parameter r_0 computed for these β values in reference [22].

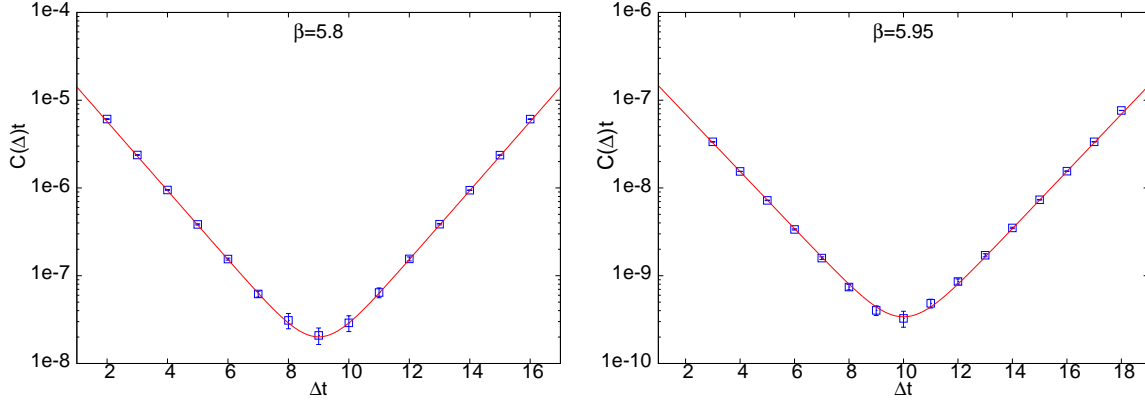
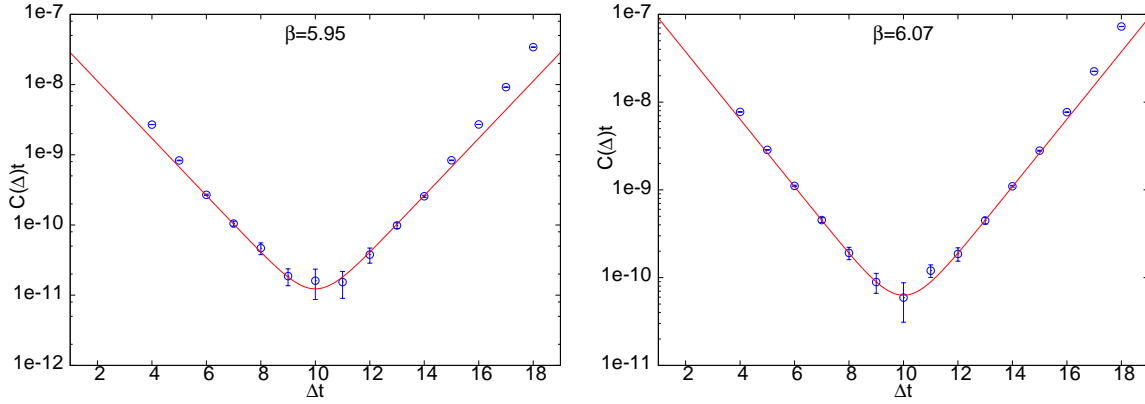
3. Results

In our simulations we could follow the glueball correlators at least upto distances of about one fermi in the scalar channel and about 0.8 fermi in the tensor channel.

To extract masses from the correlators we fitted them to the form

$$C(\Delta t) = A \left(e^{-m\Delta t} + e^{-m(T-\Delta t)} \right) \quad (3.1)$$

where m is the glueball mass and T is the full temporal extent of the lattice. Since the correlator is symmetric about $T/2$, we fold the data and use only one half of the temporal range for the fits. In figures 2 and 3 we show the scalar and tensor correlators along with the fitted correlators.

**Figure 2:** Scalar correlators at $\beta = 5.8$ (left) and 5.95 (right)**Figure 3:** Tensor correlators at $\beta = 5.95$ (left) and 6.07 (right)

Lattice	β	fit-range	ma	$\chi^2/d.o.f$
$10^3 \times 18$	5.7	5-9	0.952(11)	0.066
$12^3 \times 18$	5.8	6-9	0.906(8)	0.03
$16^3 \times 20$	5.95	5-10	0.7510(15)	0.02

Table 3: Global masses and fit parameters for the scalar channel

In addition to the masses, we also compute the effective masses from the correlators as

$$am_{\text{eff}} = -\log \frac{\langle C(\Delta t + 1) \rangle}{\langle C(\Delta t) \rangle} \quad (3.2)$$

where a is the lattice spacing. The effective masses are expected to decrease with increasing Δt and finally settle to a stable value. This stable value is an estimate of the lightest mass in the concerned J^{PC} channel. In figures 4 and 5 we plot the effective masses against Δt in the scalar and tensor channels along with masses obtained from the correlator fits (blue line). We see that at the larger values of Δt , the effective masses match with the masses from the correlators.

As a cross-check of our results, we compare our data with that reported in reference [10]. Our data is fully consistent with the results reported there and slightly more accurate.

To get an idea of the advantage of the current algorithm over the naive update algorithm, we

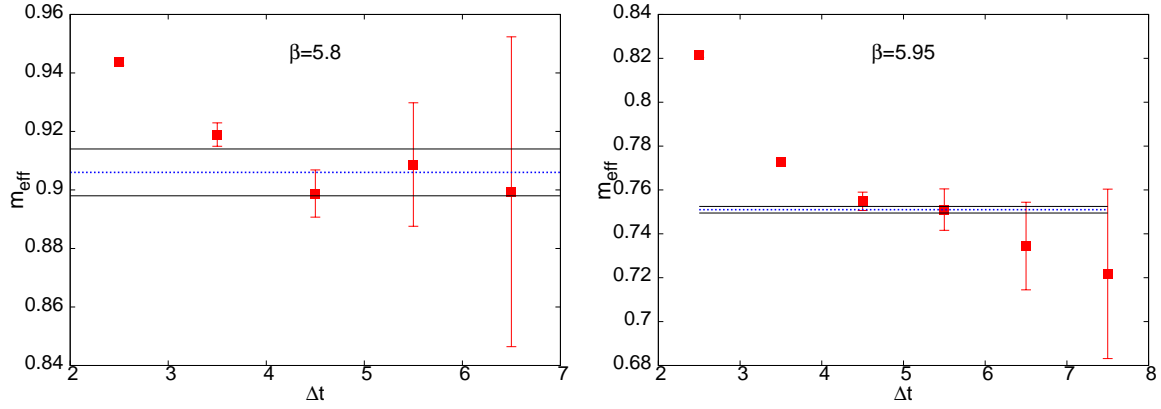


Figure 4: Effective masses in the scalar channel at β 5.8 (left) and 5.95 (right). The blue line is the mass obtained from the correlator fits. The two black lines show the error on the fitted mass.

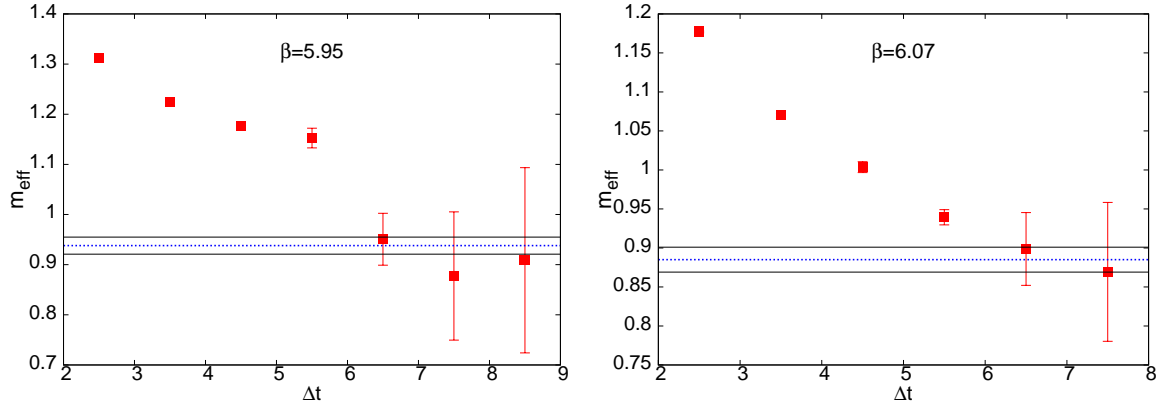


Figure 5: Effective masses in the tensor channel at β 5.95 (left) and 6.07 (right). Again the blue line is the mass from the correlator fits and the black lines show the error on the fitted mass.

Lattice	β	fit-range	ma	$\chi^2/d.o.f$
$12^3 \times 18$	5.8	4-7	1.585(54)	1.64
$12^3 \times 20$	5.95	6-10	0.938(17)	0.12
$12^3 \times 20$	6.07	6-10	0.885(16)	1.6

Table 4: Global masses and fit parameters for the tensor channel

did a few runs for the same computer time using both the methods. For these runs we used the same β values as our main run but smaller lattices. The results are reported in tables 5 and 6. We see that depending on the channel and β , the gain in time is significant and can vary from 30 to more than 700. For the tensor channel at $\beta = 6.07$ where we used a 6×6 Wilson loop, we did not manage to get a signal using the naive method. Therefore we could not calculate the gain at that point.

4. Discussions

The multilevel algorithm is very efficient for calculating quantities with very small expectation values. Operators in the tensor channel have zero expectation values and are therefore ideal for

Lattice	β	sub-lattice	iupd	loop size	run-time (mins)	$\frac{\text{error}_{naive}}{\text{error}_{multilevel}}$	gain (time)
$10^3 \times 18$	5.7	3	30	2×2	3850	5.7	32
$6^3 \times 18$	5.8	3	25	3×3	1000	5.5	30
$8^3 \times 24$	5.95	4	50	5×5	1100	18	324

Table 5: Performance comparison for Scalar Channel

Lattice	β	sub-lattice	iupd	loop size	run-time (mins)	$\frac{\text{error}_{naive}}{\text{error}_{multilevel}}$	gain (time)
$6^3 \times 18$	5.8	3	50	3×3	12000	27	729
$8^3 \times 30$	5.95	5	100	5×5	5775	20	400
$10^3 \times 30$	6.07	6	130	6×6	15000	-	-

Table 6: Performance comparison for tensor channel

direct evaluation. For scalar operators we have subtracted the non-zero vacuum expectation values from the operators to get the connected correlators directly.

Correlation functions between large loops have the advantage that they have much less contamination from excited states compared to those between elementary plaquettes. Multilevel schemes allow us to estimate the expectation values of the large loops with very high precision.

The efficiency of the algorithm depends crucially on choosing the optimal parameters for the algorithm such as the sub-lattice thickness and updates. These depend on β quite strongly. In the range of β we explored it seems that 0.5 fermi seems to be close to optimal for both the loop size and the thickness of the sub-lattice.

We observe that this error reduction technique works quite well at least in pure gauge theories. For a given computational cost, the improvement in the signal to noise ratio is several times to even a couple of orders of magnitude.

Finite volume effects is the largest source of systematic errors and to avoid them we choose our lattices such that $mL > 9$ [23]. For a more detailed discussion we refer the reader to reference [19].

Acknowledgments

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